Amendments to the Claims

This listing of claims will replace all prior versions, and listings of claims in the application.

1. (currently amended) Compounds A compound of the formula (I)

in which

- X represents C2-C4-alkyl,
- Y represents halogen, and
- Z represents C1-C4-alkyl,
- A represents alkyl,
- G represents hydrogen (a), or represents

$$\bigcap_{\mathsf{R}^1 \quad (b),} \qquad \bigcup_{\mathsf{M}}^\mathsf{L} \bigcap_{\mathsf{R}^2 \quad (c),} \qquad \bigcap_{\mathsf{SO}_{\overline{2}}} \bigcap_{\mathsf{R}^3 \quad (d),}$$

$$\bigcap_{\mathsf{R}^4 \quad \cdot} \bigcap_{\mathsf{R}^5 \quad (e),} \qquad \qquad \bigcap_{\mathsf{R}^7 \quad (g),} \bigcap_{\mathsf{$$

in which

æ

E represents a metal ion equivalent or an ammonium ion,

L represents oxygen or sulphur,

M represents oxygen or sulphur,

- R¹ represents in each case optionally substituted alkyl, alkenyl, alkoxyalkyl, alkylthioalkyl or polyalkoxyalkyl, or represents cycloalkyl or heterocyclyl, each of which is optionally substituted by halogen, alkyl or alkoxy, or represents in each case optionally substituted phenyl, heteroaryl heteroaryl, phenyl-C₁-C₄-alkyl, phenyl-C₁-C₂-alkenyl or heteryl-C₁-C₄-alkyl heteroaryl-C₁-C₄-alkyl,
- R² represents in each case optionally halogen-substituted alkyl, alkenyl, alkoxyalkyl or polyalkoxyalkyl, or represents in each case optionally substituted cycloalkyl, phenyl or benzyl,
- R³, R⁴ and R⁵ independently of one another represent in each case optionally halogen-substituted alkyl, alkoxy, alkylamino, dialkylamino, alkylthio, alkenylthio or cycloalkylthio, or represent

in each case optionally substituted phenyl, benzyl, phenoxy or phenylthio,

 R^6 and R^7 independently of one another represent hydrogen, represent in each case optionally halogen-substituted alkyl, cycloalkyl, alkenyl, alkoxy, or alkoxyalkyl, or represent in each case optionally substituted phenyl or benzyl, or R^6 and R^7 together with the N atom to which they are attached form an optionally substituted cycle which optionally contains oxygen or sulphur.

- X represents ethyl, n-propyl or n-butyl,
- Y represents halogen,
- Z represents methyl, ethyl or n-propyl,
- A represents C1-C6-alkyl,
- G represents hydrogen (a), or-represents one of the groups

$$\bigcap_{R^1 \text{ (b)}, } \bigcup_{M}^L R^2 \text{ (c)}, \qquad SO_{\frac{1}{2}} R^3 \text{ (d)},$$

$$- \bigcap_{R^5 \text{ (e)}, } R^5 \text{ (e)}, \qquad E \text{ (f)} \qquad \text{or} \qquad \bigcap_{R^7 \text{ (g)}, } R^7 \text{ (g)},$$

in which

- E represents a metal ion equivalent or an ammonium ion,
- L represents oxygen or sulphur, and
- M represents oxygen or sulphur,
- R¹ represents C₁-C₂₀-alkyl, C₂-C₂₀-alkenyl, C₁-C₆-alkoxy-C₁-C₆-alkyl, C₁-C₆-alkyl, C₁-C₆-alkyl or poly-C₁-C₄-alkoxy-C₁-C₄-alkyl, each of which is optionally mono- to heptasubstituted by halogen, mono- or disubstituted by cyano, monosubstituted by -CO-R¹¹, -C=N-OR¹¹, -CO₂R¹¹ or

CO-N $\stackrel{\mbox{R}^{11}}{\mbox{R}^{11'}}$, or represents C3-C8-cycloalkyl which is optionally monoto

trisubstituted by substituted with one, two or three substituents selected from the group consisting of halogen, C₁-C₄-alkyl [[or]] and C₁-C₄-alkoxy, wherein and in which optionally one or two not directly adjacent methylene groups of said C₃-C₈-cycloalkyl are optionally replaced by oxygen and/or or sulphur, or

represents phenyl, phenyl-C₁-C₂-alkyl or phenyl-C₁-C₂-alkenyl, each of which is optionally mono—to-trisubstituted by substituted with one, two or three substituents selected from the group consisting of halogen, cyano, nitro, C₁-C₆-alkyl, C₁-C₆-alkoxy, C₁-C₆-haloalkyl, C₁-C₆-haloalkoxy, C₁-C₆-alkylthio, C₁-C₆-alkylsulphinyl [[or]] and C₁-C₆-alkylsulphonyl, or

represents 5- or 6-membered hetaryl heteroaryl which is optionally monoor-disubstituted-by substituted with one or two substituents selected from the group consisting of halogen [[or]] and C₁-C₆-alkyl and has one or two heteroatoms selected from the group consisting of oxygen, sulphur and nitrogen,

 $R^2 \qquad \text{represents} \quad C_1-C_{20}\text{-alkyl}, \quad C_2-C_{20}\text{-alkenyl}, \quad C_1-C_6\text{-alkoxy-}C_2-C_6\text{-alkyl} \quad \text{or} \\ \\ \text{poly-}C_1-C_6\text{-alkoxy-}C_2-C_6\text{-alkyl}, \quad \text{each of which is optionally mono- to} \\ \\ \text{trisubstituted by halogen}, \quad \\ \\ \end{array}$

represents C_3 - C_8 -cycloalkyl which is optionally mono or disubstituted by substituted with one or two substituents selected from the group consisting of halogen, C_1 - C_6 -alkyl [[or]] and C_1 - C_6 -alkoxy, or

represents phenyl or benzyl, each of which is optionally mono—to trisubstituted by substituted with one, two or three substituents selected from the group consisting of halogen, cyano, nitro, C₁-C₆-alkyl, C₁-C₆-alkoxy, C₁-C₆-haloalkyl [[or]] and C₁-C₆-haloalkoxy.

- R³ represents C₁-C₈-alkyl which is optionally monor or polysubstituted by halogen, or represents phenyl or benzyl, each of which is optionally monor or disubstituted by substituted with one or two substituents selected from the group consisting of halogen, C₁-C₆-alkyl, C₁-C₆-alkoxy, C₁-C₄-haloalkyl, C₁-C₄-haloalkoxy, cyano [[or]] and nitro,
- R^4 and R^5 independently of one another represent C_1 - C_8 -alkyl, C_1 - C_8 -alkyxy, C_1 - C_8 -alkylamino, di- $(C_1$ - C_8 -alkyl)amino, C_1 - C_8 -alkylthio or C_2 - C_8 -alkenylthio, each of which is optionally mono- to trisubstituted by

halogen, or represent phenyl, phenoxy or phenylthio, each of which is optionally mone—to trisubstituted by substituted with one, two or three substituents selected from the group consisting of halogen, nitro, cyano, C₁-C₄-alkoxy, C₁-C₄-haloalkoxy, C₁-C₄-alkylthio, C₁-C₄-haloalkylthio, C₁-C₄-alkyl [[or]] and C₁-C₄-haloalkyl,

R⁶ and R⁷ independently of one another represent hydrogen, represent C₁-C₈-alkyl, C₃-C₈-cycloalkyl, C₁-C₈-alkoxy, C₃-C₈-alkenyl or C₁-C₈-alkoxy-C₂-C₈-alkyl, wherein said C₁-C₈-alkyl, C₃-C₈-cycloalkyl, C₁-C₈-alkoxy, C₃-C₈-alkenyl or C₁-C₈-alkoxy-C₂-C₈-alkyl each of which is optionally mono- to trisubstituted by halogen, or represent phenyl or benzyl, each of which is optionally mono to trisubstituted by substituted with one, two or three substituents selected from the group consisting of halogen, C₁-C₈-alkyl, C₁-C₈-haloalkyl [[or]] and C₁-C₈-alkoxy, or R⁶ and R⁷ together represent a C₃-C₆-alkylene radical which is optionally mono- or disubstituted by C₁-C₄-alkyl and in which optionally one methylene group is replaced by oxygen or sulphur,

R¹¹ represents hydrogen, or represents C₁-C₆-alkyl, C₃-C₆-alkenyl, C₃-C₆-alkynyl or C₁-C₄-alkoxy-C₂-C₄-alkyl, each of which is optionally monoto trisubstituted by halogen, or represents C₃-C₆-cycloalkyl which is optionally monoor-disubstituted by substituted with one or two substituents selected from the group consisting of halogen, C₁-C₂-alkyl [[or]] and C₁-C₂-alkoxy and in which optionally one or two not directly

adjacent methylene groups are <u>optionally</u> replaced by oxygen, or represents phenyl or phenyl- C_1 - C_3 -alkyl, each of which is optionally mono or <u>disubstituted by substituted with one or two substituents</u> selected from the group <u>consisting of halogen</u>, C_1 - C_4 -alkyl, C_1 - C_4 -alkoxy, C_1 - C_4 -haloalkyl, C_1 - C_4 -haloalkoxy, cyano [[or]] <u>and nitro, and</u>

R^{11'} represents hydrogen, C₁-C₆-alkyl or C₃-C₆-alkenyl.

- 3. (currently amended) Compounds of the formula (I) The compound according to Claim 1, in which
 - X represents ethyl or n-propyl,
 - Y represents chlorine or bromine,
 - Z represents methyl or ethyl,
 - A represents methyl, ethyl, n-propyl, n-butyl or isobutyl,
 - G represents hydrogen (a), or represents one of the groups

in which

E represents a metal ion equivalent or an ammonium ion,

- L represents oxygen or sulphur, and
- M represents oxygen or sulphur,
- R¹ represents C₁-C₁₀-alkyl, C₂-C₁₀-alkenyl, C₁-C₄-alkoxy-C₁-C₂-alkyl, C₁-C₄-alkylthio-C₁-C₂-alkyl or poly-C₁-C₃-alkoxy-C₁-C₂-alkyl, each of which is optionally mone—to pentasubstituted—by substituted with one to five substituents selected from the group consisting of fluorine [[or]] and chlorine, monosubstituted by cyano, monosubstituted by -CO-R¹¹, -C=N-OR¹¹ or CO₂R¹¹, or represents C₃-C₆-cycloalkyl which—is optionally mono—or—disubstituted—by substituted with one or two substituents selected from the group consisting of fluorine, chlorine, C₁-C₂-alkyl [[or]] and C₁-C₂-alkoxy, wherein and—in—which—optionally one or two not directly adjacent methylene groups are optionally replaced by oxygen,

represents phenyl or benzyl, each of which is optionally meno—or disubstituted by substituted with one or two substituents selected from the group consisting of fluorine, chlorine, bromine, cyano, nitro, C_1 - C_4 -alkyl, C_1 - C_4 -alkylthio, C_1 - C_4 -alkylsulphonyl, C_1 - C_4 -alkylsulphinyl, C_1 - C_2 -haloalkyl [[or]] and C_1 - C_2 -haloalkoxy, or

represents pyrazolyl, thiazolyl, pyridyl, pyrimidyl, furanyl or thienyl, each of which is optionally mono-or-disubstituted by substituted with one or two substituents selected from the group consisting of fluorine, chlorine, bromine [[or]] and C1-C2-alkyl.

R² represents C₁-C₁₀-alkyl, C₂-C₁₀-alkenyl, C₁-C₄-alkoxy-C₂-C₄-alkyl or poly-C₁-C₄-alkoxy-C₂-C₄-alkyl, each of which is optionally mone—to trisubstituted by substituted with one, two or three substituents selected from the group consisting of fluorine [[or]] and chlorine.

represents C_3 - C_7 -cycloalkyl which—is optionally monosubstituted by C_1 - C_2 -alkyl or C_1 - C_2 -alkoxy or

represents phenyl or benzyl, each of which is optionally mono—or disubstituted by substituted with one or two substituents selected from the group consisting of fluorine, chlorine, bromine, cyano, nitro, C₁-C₄-alkyl, methoxy, trifluoromethyl [[or]] and trifluoromethoxy,

- R³ represents C₁-C₄-alkyl which is optionally mono—to trisubstituted by substituted with one, two or three substituents selected from the group consisting of fluorine [[or]] and chlorine, or represents phenyl or benzyl, each of which is optionally monosubstituted by fluorine, chlorine, bromine, C₁-C₄-alkyl, C₁-C₄-alkoxy, trifluoromethyl, trifluoromethoxy, cyano or nitro,
- R⁴ and R⁵ independently of one another represent C₁-C₆-alkyl, C₁-C₆-alkyxy, C₁
 C₆-alkylamino, di-(C₁-C₆-alkyl)amino, C₁-C₆-alkylthio or C₃-C₄
 alkenylthio, each of which is optionally mone—to trisubstituted by

 substituted with one, two or three substituents selected from the group

 consisting of fluorine [[or]] and chlorine, or represent phenyl, phenoxy or

 phenylthio, each of which is optionally mone—or—disubstituted—by

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substituted with one or two substituents selected from the group consisting of fluorine, chlorine, bromine, nitro, cyano, C₁-C₃-alkoxy, trifluoromethoxy, C₁-C₃-alkylthio, C₁-C₃-alkyl [[or]] and trifluoromethyl,

- R⁶ and R⁷ independently of one another represent hydrogen, represent C₁-C₆-alkyl, C₃-C₆-cycloalkyl, C₁-C₄-alkoxy, C₃-C₆-alkenyl or C₁-C₆-alkoxy-C₂-C₆-alkyl, each of which wherein said C₁-C₆-alkyl, C₃-C₆-cycloalkyl, C₁-C₄-alkoxy, C₃-C₆-alkenyl or C₁-C₆-alkoxy-C₂-C₆-alkyl is optionally mone—to trisubstituted by substituted with one, two or three substituents selected from the group consisting of fluorine [[or]] and chlorine, represent phenyl which is optionally mone—or disubstituted by substituted with one or two substituents selected from the group consisting of fluorine, chlorine, bromine, trifluoromethyl, C₁-C₄-alkyl [[or]] and C₁-C₄-alkoxy, or R⁶ and R⁷ together represent a C₅-C₆-alkylene radical which is optionally mone—or disubstituted by methyl and in which optionally one methylene group is replaced by oxygen, and
- R¹¹ represents C₁-C₄-alkyl, C₃-C₄-alkenyl, C₃-C₄-alkynyl or C₁-C₄-alkoxy-C₂-C₃-alkyl, or represents C₃-C₆-cycloalkyl in which optionally one methylene group is replaced by oxygen.
- (currently amended) Compounds of the formula (1) The compound according to
 Claim 1 in which
 - X represents ethyl or n-propyl,

- Y represents chlorine or bromine,
- Z represents methyl or ethyl,
- A represents methyl, ethyl or n-propyl,
- G represents hydrogen (a), or represents one of the groups

in which

L represents oxygen, and

M represents oxygen or sulphur,

R¹ represents C₁-C₆-alkyl, C₂-C₆-alkenyl, C₁-C₂-alkoxy-C₁-C₂-alkyl, C₁-C₂-alkyl or poly-C₁-C₂-alkoxy-C₁-C₂-alkyl, each of which is optionally mono—to-trisubstituted by substituted with one, two or three substituents selected from the group consisting of fluorine [[or]] and chlorine, or represents cyclopropyl, cyclopentyl or cyclohexyl, each of which is optionally monosubstituted by fluorine, chlorine, methyl, ethyl or methoxy.

represents phenyl which is optionally monosubstituted by fluorine, chlorine, bromine, cyano, nitro, methyl, ethyl, n-propyl, isopropyl,

methoxy, ethoxy, methylthio, ethylthio, methylsulphinyl, ethylsulphinyl, methylsulphonyl, ethylsulphonyl, trifluoromethyl or trifluoromethoxy,

represents furanyl, thienyl or pyridyl, each of which is optionally monosubstituted by chlorine, bromine or methyl,

R² represents C₁-C₈-alkyl, C₂-C₆-alkenyl, [[or]] C₁-C₃-alkoxy-C₂-C₃-alkyl, cyclopentyl or cyclohexyl,

or represents phenyl or benzyl, each of which is optionally monosubstituted by fluorine, chlorine, bromine, cyano, nitro, methyl, methoxy, trifluoromethyl or trifluoromethoxy,

R⁶ represents hydrogen, represents C₁-C₄-alkyl, C₃-C₆-cycloalkyl, [[or]] allyl, or represents phenyl, wherein said phenyl which is optionally monosubstituted by fluorine, chlorine, bromine, methyl, methoxy or trifluoromethyl,

R⁷ represents methyl, ethyl, n-propyl, isopropyl or allyl,

 R^6 and R^7 together represent a C_3 - C_6 -alkylene radical in which optionally one methylene group is replaced by oxygen.

5. (currently amended) Process A process for preparing a compound compounds of the formula (I) according to Claim 1, characterized in that, to obtain comprising

A) compounds of the formula (I-a)

in which

A, X, Y and Z are as defined above.

condensing intramolecularly a compound compounds of the formula (II),

in which

A, X, Y and Z are as defined above in Claim 1

and

R8 represents alkyl,

are condensed intramolecularly in the presence of a diluent and in the presence of a base, to obtain a compound of the formula (I-a),

AO
$$OH_Z$$
 $(I-a)$

wherein A, X, Y and Z are as defined in Claim 1,

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(B) compounds of the formula (I b) shown above in which Λ, R[†], X, Y and Z are as defined above, compounds of the formula (I a) shown above in which Λ, X, Y and Z are as defined above are reacted

reacting a compound of the formula (I-a)

$$AO \bigvee_{OH_{Z}} O X \qquad (I-a)$$

wherein A, X, Y and Z are as defined in Claim 1,

a) with an acid halide halides of the formula (III),

in which

R¹ is as defined above in Claim 1 and

Hal represents halogen

or

B) with a carboxylic anhydride anhydrides of the formula (IV),

in which

R1 is as defined above in Claim 1,

if appropriate optionally in the presence of a diluent and if appropriate optionally in the presence of an acid binder, to obtain a compound of the formula (I-b)

AO
$$Z$$
 $(I-b)$

wherein A, R1, X, Y and Z are as defined in Claim 1,

(C) compounds of the formula (I c) shown above in which A, R², M, X, Y and

Z are as defined above and L represents oxygen, compounds of the
formula (I a) shown above in which A, X, Y and Z are as defined above
are in each case reacted

reacting a compound of the formula (I-a)

wherein A, X, Y and Z are as defined in Claim 1,

with a chloroformic ester esters or a chloroformic thioester thioesters of the formula (V).

in which

R² and M are as defined above in Claim 1,

if appropriate optionally in the presence of a diluent and if appropriate optionally in the presence of an acid binder, to obtain a compound of the formula (I-c)

wherein A, R², M, X, Y and Z are as defined in Claim 1, and L is oxygen,

(D) compounds of the formula (I-c) shown above in which A, R², M, X, Y and Z-are as defined above and L represents sulphur, compounds of the formula (I a) shown above in which A, X, Y and Z are as defined above are in each case reacted

reacting a compound of the formula (I-a)

$$AO \bigvee_{OH_{Z}} \bigvee_{Y} \underbrace{(I-a)}_{Y}$$

wherein A, X, Y and Z are as defined in Claim 1.

 α) with <u>a</u> chloromonothioformic <u>ester</u> esters or <u>a</u> chlorodithioformic <u>ester</u> esters of the formula (VI),

in which

M and R² are as defined above in Claim 1,

if appropriate optionally in the presence of a diluent and if appropriate optionally in the presence of an acid binder,

or

 β) with carbon disulphide and then with <u>a compound</u> eompounds of the formula (VII),

in which

R2 is as defined above in Claim 1 and

Hal represents chlorine, bromine or iodine,

if appropriate optionally in the presence of a diluent and if appropriate optionally in the presence of a base, to obtain a compound of the formula (I-c)

wherein A, R², M, X, Y and Z are as defined in Claim 1, and L is sulphur,

(E) compounds of the formula (I d) shown above in which A, R³, X, Y and Z are as defined above, compounds of the formula (I a) shown above in which A, X, Y and Z are as defined above are in each case reacted

wherein A, X, Y and Z are as defined in Claim 1,

reacting a compound of the formula (I-a)

with a sulphonyl chloride ehlorides of the formula (VIII),

in which

R³ is as defined above in Claim 1,

if appropriate optionally in the presence of a diluent and if appropriate optionally in the presence of an acid binder, to obtain a compound of the formula (I-d)

wherein A, R³, X, Y, and Z are as defined in Claim 1,

(F) compounds of the formula (I e) shown above in which Λ, L, R⁴, R⁵, X, Y and Z are as defined above, compounds of the formula (I a) shown above in which Λ, X, Y and Z are as defined above are in each case reacted reacting a compound of the formula (I-a)

wherein A, X, Y and Z are as defined in Claim 1,

with a phosphorus compound compounds of the formula (IX),

in which

L, R⁴ and R⁵ are as defined above in Claim 1 and

Hal represents halogen,

if appropriate optionally in the presence of a diluent and if appropriate optionally in the presence of an acid binder, to obtain a compound of the formula (I-e)

wherein A, L, R4, R5, X, Y and Z are as defined in Claim 1,

(G) compounds of the formula (I-f) shown above in which A, E, X, Y and Z are as defined above, compounds of the formula (I a) in which A, X, Y and Z are as defined above are in each case reacted

reacting a compound of the formula (I-a)

AO
$$OH_Z$$
 $(I-a)$

wherein A, X, Y and Z are as defined in Claim 1,

with <u>a</u> metal <u>compound</u> eompounds or <u>amine</u> amines of the formulae (X) or (XI), respectively,

$$R^{10}$$
, R^{9}

Me(OR⁸), (X)

 R^{10} , R^{9}

(XI)

in which

Me represents a mono- or divalent metal,

t represents the number 1 or 2 and

R⁸, R⁹, and R¹⁰ independently of one another represent hydrogen or alkyl,

if appropriate optionally in the presence of a diluent, to obtain a compound of the formula (I-f)

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wherein A, E, X, Y and Z are as defined in Claim 1,

(H) compounds of the formula (I g) shown above in which A, L, R⁶, R⁷, X, Y and Z are as defined above, compounds of the formula (I a) shown above in which A, X, Y and Z are as defined above are in each case reacted reacting a compound of the formula (I-a)

AO
$$OH_Z$$
 $(I-a)$

wherein A, X, Y and Z are as defined in Claim 1,

 α) with <u>an isocyanate</u> isocyanates or <u>an isothiocyanate</u> isothiocyanates of the formula (XII).

$$R^6$$
-N=C=L (XII)

in which

R⁶ and L are as defined above in Claim 1,

if appropriate optionally in the presence of a diluent and if appropriate optionally in the presence of a catalyst, or

 B) with a carbamoyl <u>chloride</u> ehlorides or a thiocarbamoyl <u>chloride</u> ehlorides of the formula (XIII).

$$R^6 \longrightarrow N$$
 CI (XIII)

in which

L, R⁶ and R⁷ are as defined above in Claim 1,

if appropriate optionally in the presence of a diluent and if appropriate optionally in the presence of an acid binder, to obtain a compound of the formula (I-g)

AO
$$\frac{1}{R^7-N}$$
 $\frac{(I-g)}{R^6}$

wherein A, L, R6, R7, X, Y and Z are as defined in Claim 1,

(I) eempounds of the formulae (I-a) to (I-g) shown above, cis/trans isomer mixtures—of—formulae—(I-a') to (I-g'), known, for example, from EP A-835-243.

separating the cis/trans isomer mixture of a compound of formula (I-a')-(I-g'):

$$A-O \longrightarrow \begin{array}{c} H & O \\ O & Z \end{array}$$

$$(I-b)_{i}$$

$$A-O \longrightarrow \begin{array}{c} H \\ \hline \\ L \\ \hline \\ M \\ R^2 \end{array} \qquad (I-c')$$

$$A-OW \xrightarrow{H} V X Y (I-d')_{i}$$

$$A-OW \bigvee_{E-O} \bigvee_{Z} \bigvee_{Y} (I-f), or$$

$$A-O \longrightarrow \begin{array}{c} H \\ \hline \\ R^7-N \\ \hline \\ D^6 \end{array} \qquad (I-g')$$

in which

A, E, L, M X, Y, Z, R¹, R², R³, R⁴, R⁵, R⁶ and R⁷ are as defined above in Claim 1, are separated using a physical separation process processes, such as, for example, column chromatography or fractional crystallization, to obtain the respective cis-isomer of the formula (I-a)-(I-g):

$$\begin{array}{c|c} & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & & \\ & &$$

AO

R²-M

O

Z

(I-e),

AO

R³-SO₂-O

Z

(I-e),

AO

R⁵- |

AO

R⁵- |

AO

$$R^5$$
- |

 R^5 - |

(J) eompounds of the formula (I-a), compounds hydrolyzing a compound of the formula formulae (I-b), (I-c), (I-d), (I-e), (I-f) or (I-g);

AO
$$Z$$
 R^7-N
 R^6

in which A, E, L, M, X, Y, Z, R¹, R², R³, R⁴, R⁵, R⁶ and R⁷ are as defined above in Claim 1 are hydrolysed using, for example, aqueous bases and then acidified acidifying to obtain a compound of formula (I-a)

$$AO \underbrace{\hspace{1cm} \bigcup_{OH_{Z}}^{O} X}_{OH_{Z}} \underbrace{\hspace{1cm} (I-a)}_{I}$$

wherein A, X, Y and Z are as defined in Claim 1.

6. (cancelled)

- 7. (currently amended) Pesticides and/or herbicides, characterized in that they comprise Δ pesticide or a herbicide preparation, comprising at least one compound of the formula
 (f) according to Claim 1.
- 8. (currently amended) <u>A method</u> <u>Method</u> for controlling animal pests and/or or unwanted vegetation, characterized in that compounds comprising contacting a compound of the formula (f) according to Claim 1 are allowed to act on with pests and/or or their habitat or unwanted vegetation.

- 9. (cancelled)
- 10. (currently amended) Process A process for preparing pesticides and/or herbicides, eharacterized in that compounds a pesticide or a herbicide preparation, comprising mixing a compound of the formula (I) according to Claim 1 are mixed with one or more extenders and/or or surfactants, or combinations thereof.
- 11. (currently amended)

 A composition Composition, comprising an effective
 amount of a combination of active compounds comprising
 - (a') at least one substituted eyelie ketoenol compound of the formula (I) in which A, G, X, Y and Z are as defined above according to Claim 1

and

(b') at least one crop plant compatibility-improving compound <u>selected</u> from the <u>following</u> group <u>consisting</u> of <u>eompounds</u>:

4-dichloroacetyl-1-oxa-4-azaspiro[4.5]decane (AD-67, MON-4660), 1-dichloroacetylhexahydro-3,3,8a-trimethylpyrrolo[1,2-a]pyrimidin-6(2H)-one (dicyclonon, BAS-145138), 4-dichloroacetyl-3,4-dihydro-3-methyl-2H-1,4-benzoxazine (benoxacor), 1-methylhexyl 5-chloroquinoline-8-oxyacetate (cloquintocet-mexyl [[-]] ef. also-related compounds in EP-A-86750, EP-A-94349, EP-A-191736, EP-A-492366), 3-(2-chlorobenzyl)-1-(1-methyl-1-phenylethyl)urea (cumyluron), α -(cyanomethoximino)phenylacetonitrile (cyometrinil), 2,4-dichlorophenoxyacetic acid (2,4-D), 4-(2,4-dichlorophenoxy)butyric acid (2,4-DB), 1-(1-methyl-1-phenylethyl)-3-(4-methylphenyl)urea (daimuron, dymron), 3,6-dichloro-

2-methoxybenzoic acid (dicamba), S-1-methyl 1-phenylethyl piperidine-1-thiocarboxylate (dimeninerate). 2.2-dichloro-N-(2-oxo-2-(2-propenylamino)ethyl)-N-(2-propenyl)acetamide (DKA-24). 2.2-dichloro-N.N-di-2propenylacetamide (dichlormid), 4,6-dichloro-2-phenylpyrimidine (fenclorim), ethyl 1-(2,4-dichlorophenyl)-5-trichloromethyl-1H-1,2,4-triazole-3-carboxylate (fenchlorazole-ethyl [[-]] ef. also related compounds in EP A-174562 and EP A-346620). phenylmethyl 2-chloro-4-trifluoromethylthiazole-5-carboxylate (flurazole). 4-chloro-N-(1,3-dioxolan-2-ylmethoxy)-α-trifluoroacetophenone 3-dichloroacetyl-5-(2-furanyl)-2,2-dimethyloxazolidine oxime (fluxofenim). (furilazole, MON-13900), ethyl 4,5-dihydro-5,5-diphenyl-3-isoxazolecarboxylate (isoxadifen-ethyl [[-]] ef. also related compounds in WO A 95/07897). 1-(ethoxycarbonyl)ethyl 3,6-dichloro-2-methoxybenzoate (lactidichlor), (4-chloroo-tolyloxy)acetic acid (MCPA), 2-(4-chloro-o-tolyloxy)propionic acid (mecoprop), diethyl 1-(2,4-dichorophenyl)-4,5-dihydro-5-methyl-1H-pyrazole-3,5-dicarboxylate (mefenpyr-diethyl [[-]] ef. also related compounds in WO-A-91/07874), 2-dichloromethyl-2-methyl-1,3-dioxolane (MG-191), 2-propenyl 1oxa-4-azaspiro[4.5]decane-4-carbodithioate (MG-838), 1,8-naphthalic anhydride, α-(1,3-dioxolan-2-ylmethoximino)phenylacetonitrile (oxabetrinil), 2,2-dichloro-N-(1,3-dioxolan-2-ylmethyl)-N-(2-propenyl)acetamide (PPG-1292), 3-dichloroacetyl-2,2-dimethyloxazolidine (R-28725), 3-dichloroacetyl-2,2,5-trimethyloxazolidine (R-29148), 4-(4-chloro-o-tolyl)butyric acid, 4-(4-chlorophenoxy)butyric acid. diphenylmethoxyacetic acid, methyl diphenylmethoxyacetate. ethyl diphenylmethoxyacetate, methyl 1-(2chlorophenyl)-5-phenyl-1H-pyrazole-3-carboxylate. 1_ ethvl (2,4-dichlorophenyl)-5-methyl-1H-pyrazole-3-carboxylate, ethyl 1-(2,4-dichlorophenyl)-5-isopropyl-1H-pyrazole-3-carboxylate, ethyl 1-(2.4-dichlorophenyl)-5-(1,1-dimethylethyl)-1H-pyrazole-3-carboxylate, ethyl 1-(2,4-dichlorophenyl)-5phenyl-1H-pyrazole-3-carboxylate (ef. also related compounds in EP-A-269806 and EP-A-333131), ethyl 5-(2.4-dichlorobenzyl)-2-isoxazoline-3-carboxylate. ethyl 5-phenyl-2-isoxazoline-3-carboxylate, ethyl 5-(4-fluorophenyl)-5-phenyl-2isoxazoline-3-carboxylate (cf. also related compounds in WO A 91/08202). 1,3-dimethylbut-1-vl 5-chloroguinoline-8-oxyacetate. 4-allyloxybutyl 5-chloroguinoline-8-oxyacetate. 1-allyloxyprop-2-yl 5-chloroguinoline-8oxyacetate, methyl 5-chloroquinoxaline-8-oxyacetate, ethyl 5-chloroquinoline-8ox vacetate. allvl 5-chloroquinoxaline-8-oxvacetate. 2-oxoprop-1-vl 5-chloroquinoline-8-oxyacetate, diethyl 5-chloroquinoline-8-oxymalonate, diallyl 5-chloroquinoxaline-8-oxymalonate, diethyl 5-chloroquinoline-8-oxymalonate (cf. also related compounds in EP A 582198), 4-carboxychroman-4-ylacetic acid (AC-304415, cf. EP-A-613618), 4-chlorophenoxyacetic acid, 3,3'-dimethyl-4methoxybenzophenone, 1-bromo-4-chloromethylsulphonylbenzene, 1-[4-(N-2methoxybenzoylsulphamoyl)phenyl]-3-methylurea (also known N-(2methoxybenzoyl)-4-[(methylaminocarbonyl)amino]benzenesulphonamide), 1-[4-(N-2-methoxybenzoylsulphamoyl)phenyl]-3,3-dimethylurea, 1-[4-(N-4.5dimethylbenzoylsulphamoyl)phenyl]-3-methylurea,

1-[4-(N-naphthylsulphamoyl)phenyl]-3,3-dimethylurea, and N-(2-methoxy-5-

methylbenzoyl)-4-(cyclopropylaminocarbonyl)benzenesulphonamide,

and/or one of the following compounds, defined by general formulae, a compound of the general formula (IIa)

[[or]] a compound of the general formula (IIb)

[[or]] a compound of the formula (IIc)

where

m represents a number 0, 1, 2, 3, 4 or 5,

A¹ represents one of the <u>following</u> divalent heterocyclic <u>groups groupings shown</u> below;

$$R^{17}$$
 R^{18}
 R^{17}
 R^{18}
 R^{19}
 R^{18}
 R^{19}

- n represents a number 0, 1, 2, 3, 4 or 5,
- A² represents optionally C₁-C₄-alkyl and/or C₁-C₄-alkoxyearbonyl-substituted alkanediyl having 1 or 2 carbon atoms optionally substituted with one or more substituents selected from the group consisting of C₁-C₄-alkyl and C₁-C₄-alkoxyearbonyl,
- R¹² represents hydroxy! hydroxy, mercapto, amino, C₁-C₇-alkoxy, C₁-C₆-alkenyloxy, C₁-C₆-alkenyloxy-C₁-C₆-alkoxy, C₁-C₆-alkylthio, C₁-C₆-alkylamino or di-(C₁-C₄-alkyl)amino,
- R¹³ represents hydroxyl hydroxy, mercapto, amino, C₁-C₆-alkoxy, C₁-C₆-alkylthio, C₁-C₆-alkylamino or di-(C₁-C₄-alkyl)amino,
- R¹⁴ represents in each case optionally fluorine, chlorine and/or bromine substituted
 C₁-C₄-alkyl optionally substituted with one or more substituents selected from the group consisting of fluorine, chlorine and bromine,
- R¹⁵ represents hydrogen, in each case optionally fluorine, chlorine and/or bromine-substituted C₁-C₆-alkyl, C₂-C₆-alkenyl or C₂-C₆-alkynyl in each case optionally substituted with one or more substituents selected from the group consisting of fluorine, chlorine and bromine, C₁-C₄-alkoxy-C₁-C₄-alkyl, dioxolanyl-C₁-C₄-

alkyl, furyl, furyl-C₁-C₄-alkyl, thienyl, thiazolyl, piperidinyl, [[or]] <u>phenyl</u> optionally <u>substituted with one or more substituents selected from the group consisting of fluorine[[-]], chlorine[[-]] and/or and bromine[[-]], or C₁-C₄-alkyl-substituted phenyl,</u>

- R¹⁶ represents hydrogen, in-each case optionally fluorine, chlorine and/or-brominesubstituted C₁-C₆-alkyl, C₂-C₆-alkenyl or C₂-C₆-alkynyl in each case optionally
 substituted with one or more substituents selected from the group consisting of
 fluorine, chlorine and bromine, C₁-C₄-alkoyy-C₁-C₄-alkyl, dioxolanyl-C₁-C₄alkyl, furyl, furyl-C₁-C₄-alkyl, thiazolyl, piperidinyl, [[or]] phenyl
 optionally substituted with one or more substituents selected from the group
 consisting of fluorine[[-]], chlorine[[-]] and/or and bromine[[-]], or C₁-C₄-alkylsubstituted phenyl, or R¹⁵ and R¹⁶ together also represent C₃-C₆-alkanediyl or C₂C₃-oxaalkanediyl, each of which is optionally substituted by C₁-C₄-alkyl, phenyl,
 furyl, a fused benzene ring or by two substituents which, together with the C
 atom to which they are attached, form a 5- or 6-membered carbocycle,
- R¹⁷ represents hydrogen, cyano, or halogen, or represents in each case optionally fluorine, chlorine and/or bromine substituted C₁-C₄-alkyl, C₃-C₆-cycloalkyl or phenyl in each case optionally substituted with one or more substituents selected from the group consisting of fluorine, chlorine and bromine,
- R^{18} represents hydrogen or optionally [[hydroxyl-]] hydroxy-, cyano-, halogen- or C_1 - C_4 -alkoxy-substituted C_1 - C_6 -alkyl, C_3 - C_6 -cycloalkyl or tri(C_1 - C_4 -alkyl)silyl,

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- R¹⁹ represents hydrogen, cyano, or halogen, or represents in each case optionally fluorine, ehlorine and/or bromine substituted C₁-C₄-alkyl, C₃-C₆-cycloalkyl or phenyl in each case optionally substituted with one or more substituents selected from the group consisting of fluorine, chlorine and bromine,
- X^1 represents nitro, cyano, halogen, C_1 - C_4 -alkyl, C_1 - C_4 -haloalkyl, C_1 - C_4 -alkoxy or C_1 - C_4 -haloalkoxy,
- X² represents hydrogen, cyano, nitro, halogen, C₁-C₄-alkyl, C₁-C₄-haloalkyl, C₁-C₄-alkoxy or C₁-C₄-haloalkoxy,
- X³ represents hydrogen, cyano, nitro, halogen, C₁-C₄-alkyl, C₁-C₄-haloalkyl, C₁-C₄-alkoxy or C₁-C₄-haloalkoxy,

and/or the following compounds, defined by general formulae, or selected from the group consisting of a compound of the general formula (IId)

$$O \xrightarrow{R^{21}} (X^5)_{\nu} \xrightarrow{R^{20}} (X^4)_{i}$$

$$SO_2 \xrightarrow{Q} (IId)$$

[[or]] and a compound of the general formula (IIe)

$$R^{23} \xrightarrow[R^{24}]{} SO_2^{(X^5)_v}$$
 (IIe)

where

t represents a number 0, 1, 2, 3, 4 or 5,

v represents a number 0, 1, 2, 3, 4 or 5,

 R^{20} represents hydrogen or C_1 - C_4 -alkyl,

R²¹ represents hydrogen or C₁-C₄-alkyl,

R²² represents hydrogen, in each case optionally cyano-, halogen- or C₁-C₄-alkoxy-substituted C₁-C₆-alkyl, C₁-C₆-alkoxy, C₁-C₆-alkylthio, C₁-C₆-alkylamino or di-(C₁-C₄-alkyl)amino, or in each case optionally cyano-, halogen- or C₁-C₄-alkyl-substituted C₃-C₆-cycloalkyl, C₃-C₆-cycloalkyloxy, C₃-C₆-cycloalkylthio or C₃-C₆-cycloalkylamino,

R²³ represents hydrogen, optionally cyano-, [[hydroxyl-]] <u>hydroxy-</u>, halogen- or C₁-C₄-alkoxy-substituted C₁-C₆-alkyl, in each case optionally cyano- or halogen-substituted C₃-C₆-alkenyl or C₃-C₆-alkynyl, or optionally cyano-, halogen- or C₁-C₄-alkyl-substituted C₃-C₆-cycloalkyl,

R²⁴ represents hydrogen, optionally cyano-, [[hydroxyl-]] <u>hydroxy-</u>, halogen- or C₁-C₄-alkoxy-substituted C₁-C₆-alkyl, in each case optionally cyano- or halogen-

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substituted C_3 - C_6 -alkenyl or C_3 - C_6 -alkynyl, optionally cyano-, halogen- or C_1 - C_4 -alkyl-substituted C_3 - C_6 -cycloalkyl, or optionally nitro-, cyano-, halogen-, C_1 - C_4 -alkyl-, C_1 - C_4 -haloalkyl-, C_1 - C_4 -alkoxy- or C_1 - C_4 -haloalkoxy-substituted phenyl, or R^{22} together with R^{23} represents represent in each case optionally C_1 - C_4 -alkyl-substituted C_2 - C_6 -alkanediyl or C_2 - C_5 -oxaalkanediyl,

- X⁴ represents nitro, cyano, carboxyl, carbamoyl, formyl, sulphamoyl, hydroxyl hydroxy, amino, halogen, C₁-C₄-alkyl, C₁-C₄-haloalkyl, C₁-C₄-alkoxy or C₁-C₄-haloalkoxy, and
- X⁵ represents nitro, cyano, carboxyl, carbamoyl, formyl, sulphamoyl, hydroxyl hydroxy, amino, halogen, C₁-C₄-alkyl, C₁-C₄-haloalkyl, C₁-C₄-alkoxy or C₁-C₄-haloalkoxy.
- 12. (currently amended) Composition The composition according to Claim 11, where the crop plant compatibility-improving compound is selected from the following group consisting of compounds:

cloquintocet-mexyl, fenchlorazole-ethyl, isoxadifen-ethyl, mefenpyr-diethyl, furilazole, fenclorim, cumyluron, dymron, or the compounds He 5 or He 11

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$$\bigcap_{H} \bigcap_{SO_2} \bigcap_{N} \bigcap_{CH_3} CH_3$$

- 13. (currently amended) Composition The composition according to Claim 41-or 12, where the crop plant compatibility-improving compound is cloquintocet-mexyl or mesenpyr-diethyl.
- 14. (currently amended) Method A method for controlling unwanted vegetation, eharacterized in that comprising contacting a composition according to Claim 11 is allowed to act on the plants or their habitat with unwanted vegetation.
- 15. (cancelled)
- 16. (currently amended) Compounds A compound of the formula (II)

in which

A represents alkyl,

X represents C₂-C₄-alkyl,

Y represents halogen,

Z represents C₁-C₄-alkyl, and

R⁸ represents alkyl.

A. X. Y. Z and R8 are as defined above.

17. (currently amended) Compounds A compound of the formula (XVI)

in which

A represents alkyl,

X represents C₂-C₄-alkyl,

Y represents halogen, and

Z represents C₁-C₄-alkyl.

A, X, Y and Z are as defined above.